

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050319\
 Data File : BM020117.D
 Acq On : 03 May 2019 18:16
 Operator : JU/SJ
 Sample : K2584-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSSI-0419-S-DH-7

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM043019.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.011	3	4	12	rVB	117517	104903	1.17%	0.179%
2	5.363	397	404	412	rBV	3426462	4684180	52.26%	8.011%
3	6.951	665	674	687	rBV	3119181	5320923	59.36%	9.100%
4	7.316	729	736	745	rBV	3404233	5211402	58.14%	8.913%
5	7.787	809	816	822	rBV	476019	743954	8.30%	1.272%
6	8.104	862	870	878	rBV	2420883	3825725	42.68%	6.543%
7	8.951	1006	1014	1024	rBV	1932657	3131758	34.94%	5.356%
8	10.575	1284	1290	1298	rBV	519030	878665	9.80%	1.503%
9	13.045	1703	1710	1717	rBV	4292074	6267446	69.93%	10.719%
10	13.886	1847	1853	1860	rVB	340634	482280	5.38%	0.825%
11	14.427	1939	1945	1953	rBV2	737942	1119318	12.49%	1.914%
12	15.916	2191	2198	2212	rBV	4071877	5467479	61.00%	9.351%
13	17.174	2405	2412	2416	rBV	968958	1380689	15.40%	2.361%
14	17.215	2416	2419	2425	rVB	72428	95160	1.06%	0.163%
15	18.057	2557	2562	2573	rBV	530229	823885	9.19%	1.409%
16	19.227	2757	2761	2767	rVB	232189	340876	3.80%	0.583%
17	19.339	2775	2780	2791	rBV2	466371	661020	7.37%	1.131%
18	19.592	2820	2823	2832	rVB	194724	279320	3.12%	0.478%
19	19.798	2852	2858	2864	rBV	7407061	8963087	100.00%	15.329%
20	21.056	3068	3072	3080	rBV2	406937	509508	5.68%	0.871%
21	21.356	3116	3123	3127	rBV	1340590	1890641	21.09%	3.234%
22	21.392	3127	3129	3134	rVB	125860	134994	1.51%	0.231%
23	21.498	3143	3147	3151	rBV2	200359	239524	2.67%	0.410%
24	21.939	3218	3222	3228	rBV	282990	367386	4.10%	0.628%
25	22.409	3297	3302	3308	rBV	421930	562996	6.28%	0.963%
26	22.539	3321	3324	3329	rVB	147763	198063	2.21%	0.339%
27	22.927	3385	3390	3407	rVB2	492025	1102733	12.30%	1.886%
28	23.509	3484	3489	3499	rVB2	429399	874951	9.76%	1.496%
29	23.639	3505	3511	3528	rVB	920436	1850529	20.65%	3.165%
30	24.174	3597	3602	3610	rVB2	320927	618022	6.90%	1.057%
31	24.939	3728	3732	3740	rVB3	163943	338426	3.78%	0.579%

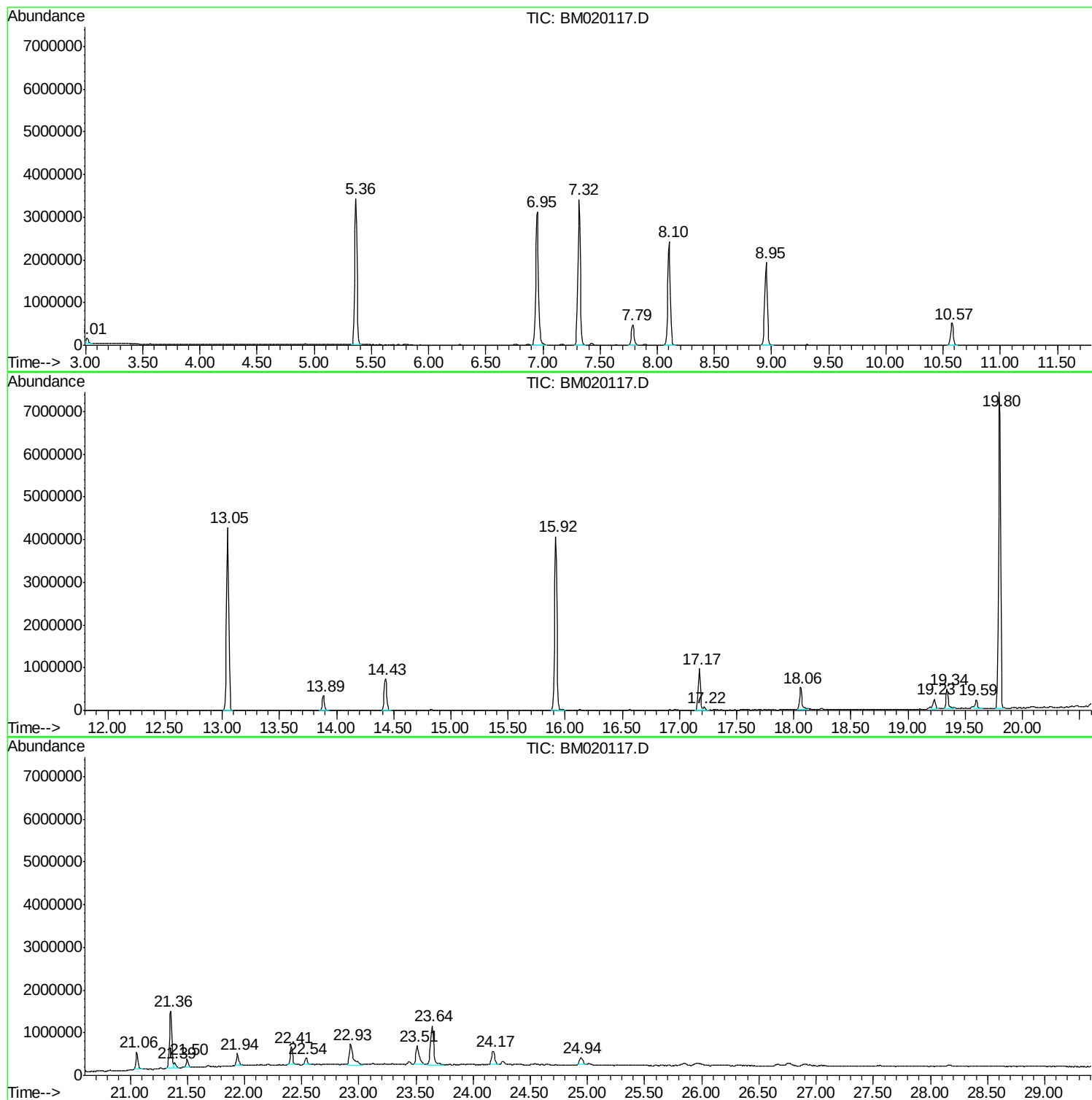
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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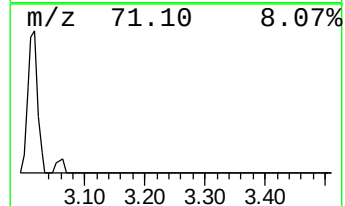
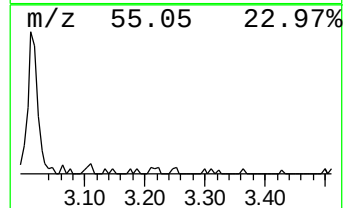
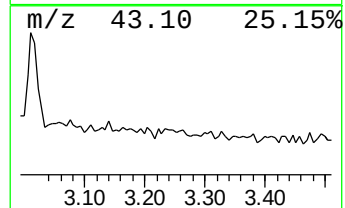
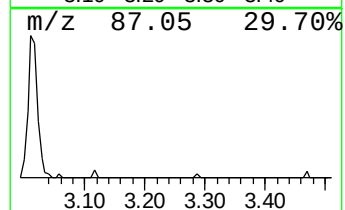
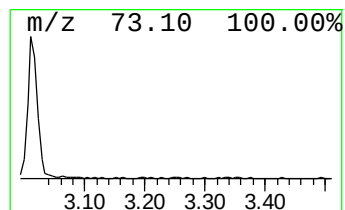
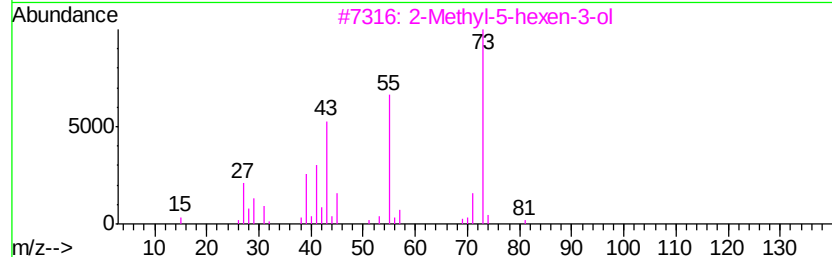
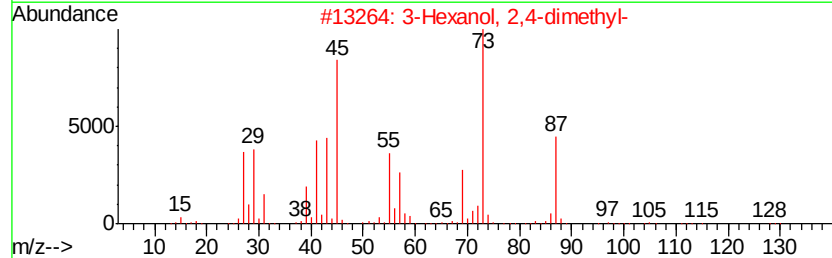
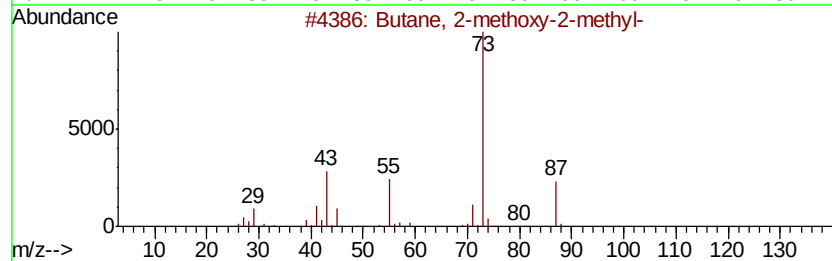
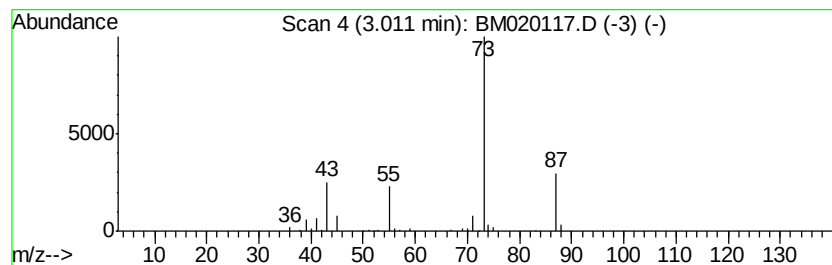
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown3.01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.01	2.82 ng	104903	1,4-Dichlorobenzene-d4	7.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	45
2			3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	9
3			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	9
4			1-Propene-1-thiol	74	C3H6S	000925-89-3	7
5			Borane, dimethoxy-	74	C2H7BO2	004542-61-4	5



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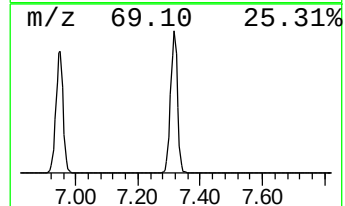
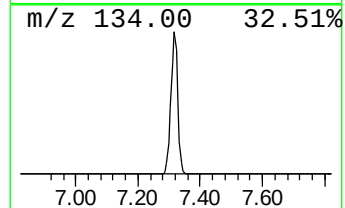
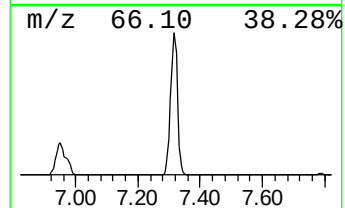
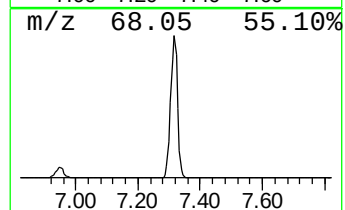
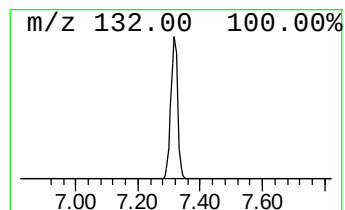
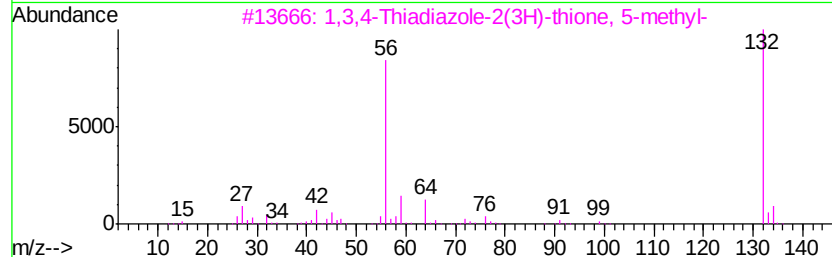
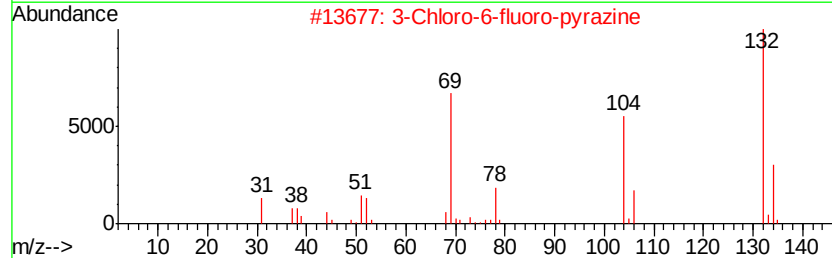
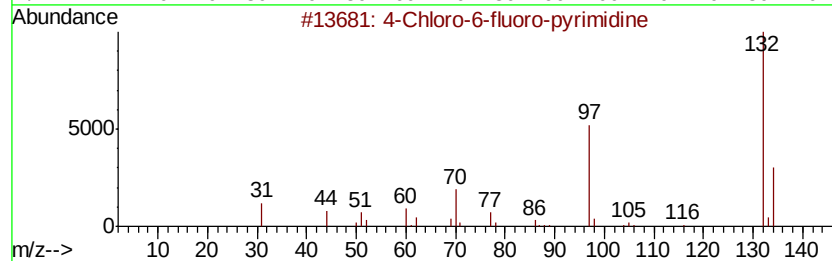
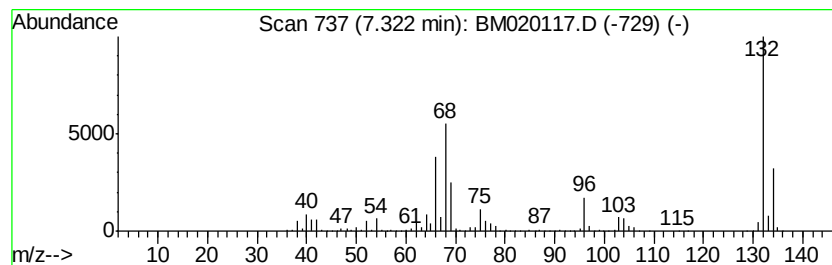
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Peak Number 2 unknown7.32 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.32	140.10 ng	5211400	1,4-Dichlorobenzene-d4	7.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
3		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	10
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	10
5		3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	9



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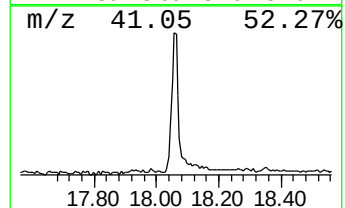
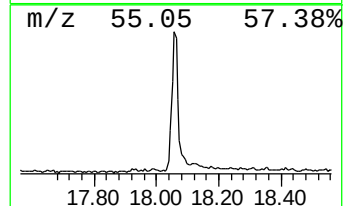
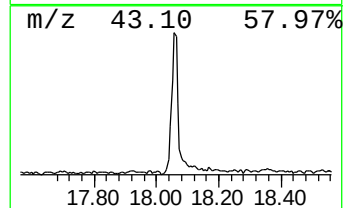
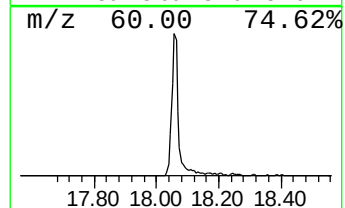
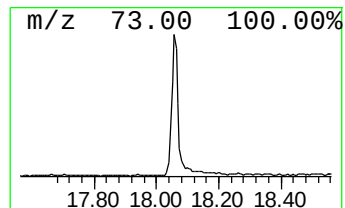
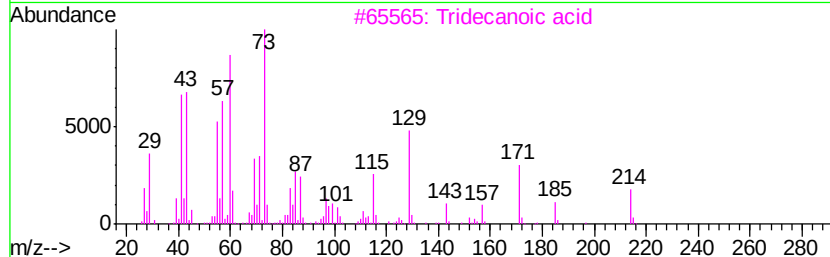
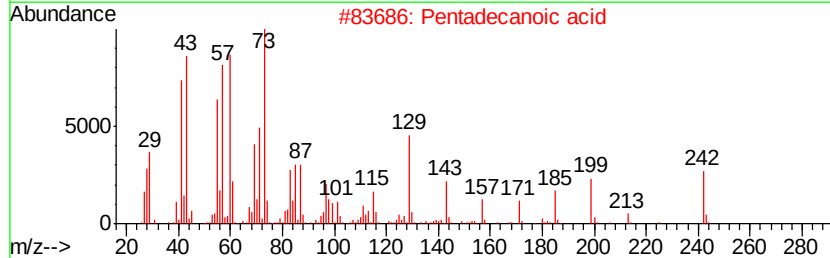
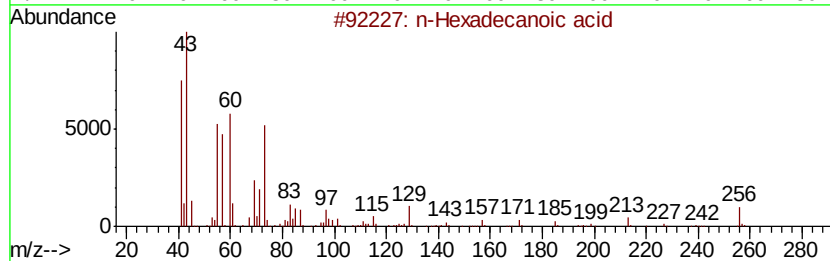
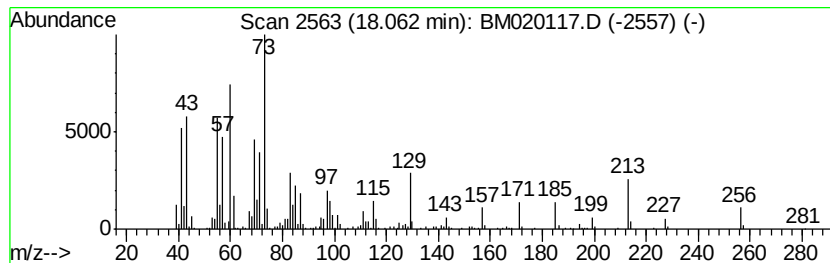
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.06	11.93 ng	823885	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	58
3		Tridecanoic acid	214	C13H26O2	000638-53-9	55
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	50
5		n-Decanoic acid	172	C10H20O2	000334-48-5	43



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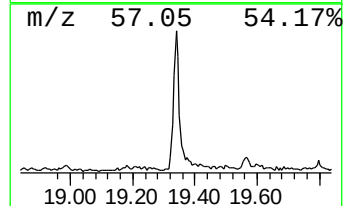
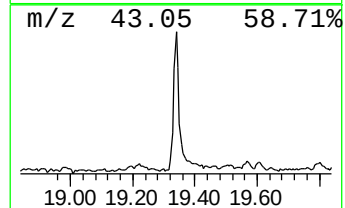
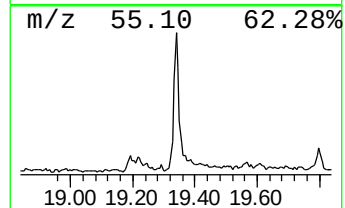
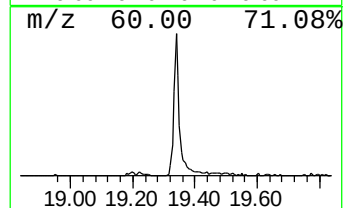
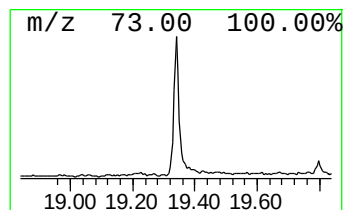
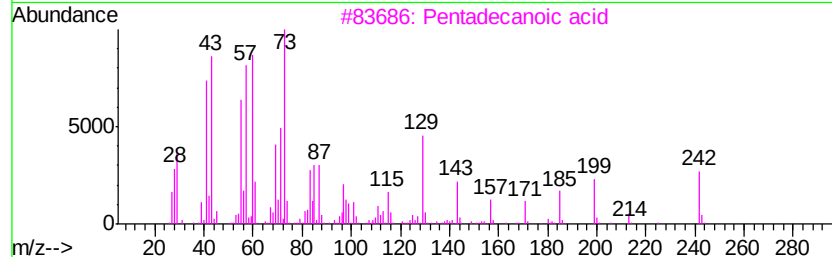
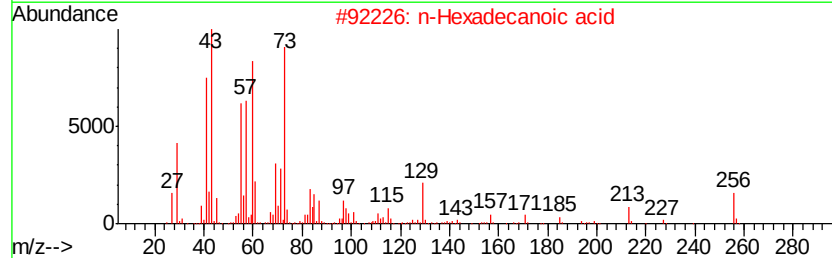
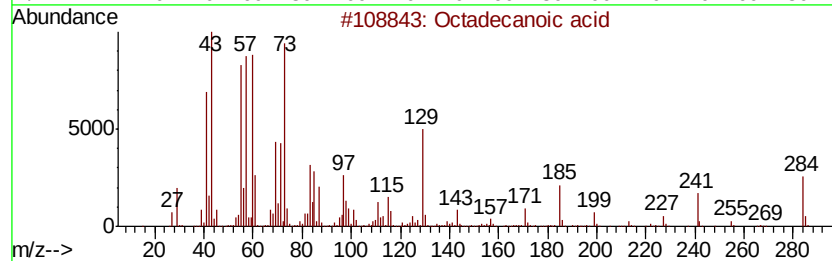
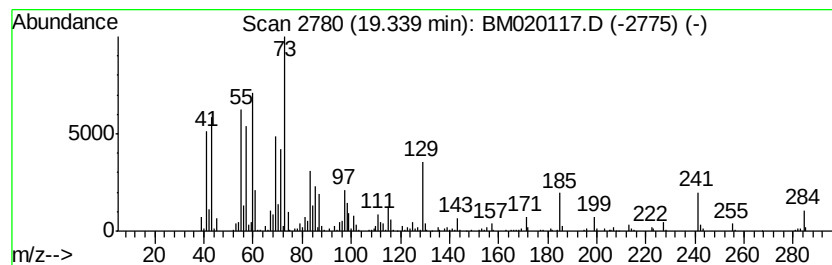
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Peak Number 5 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
19.34	6.99 ng	661020	Chrysene-d12		21.36
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	97
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
4	Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	72
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	58



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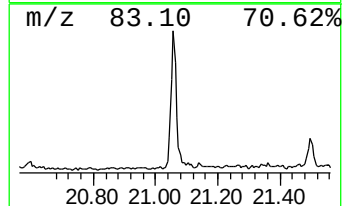
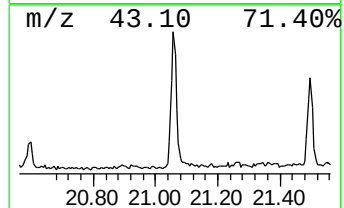
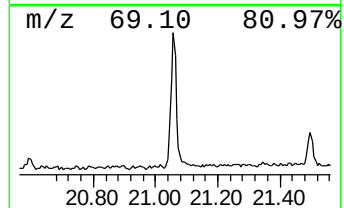
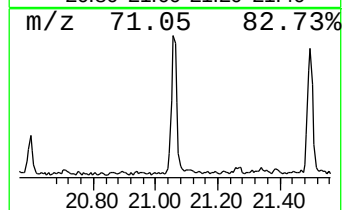
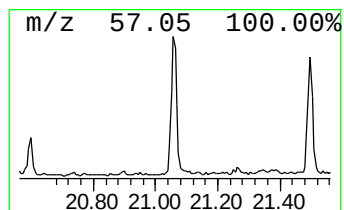
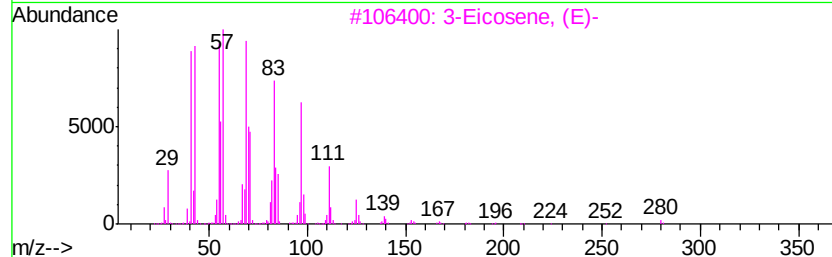
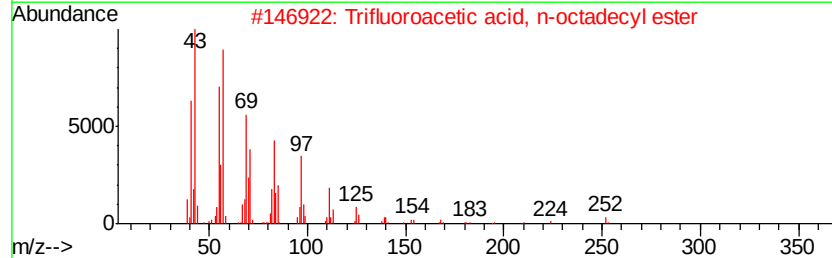
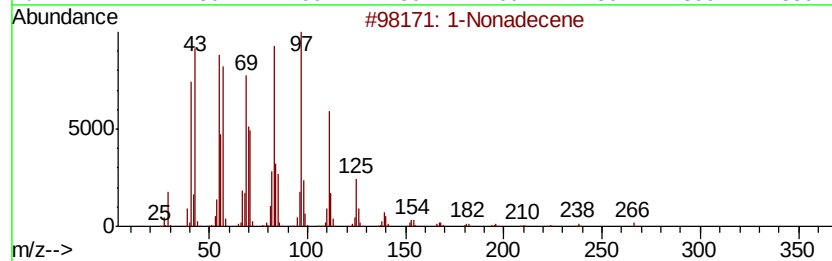
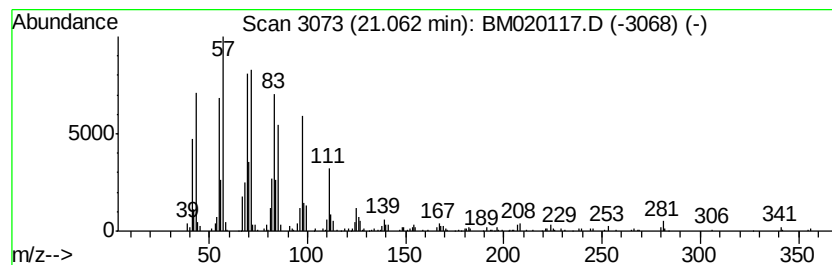
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Peak Number 6 1-Nonadecene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.06	5.39 ng	509508	Chrysene-d12	21.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonadecene	266	C19H38	018435-45-5	89
2		Trifluoroacetic acid, n-octadecyl...	366	C20H37F3O2	079392-43-1	87
3		3-Eicosene, (E)-	280	C20H40	074685-33-9	81
4		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	80
5		11-Tricosene	322	C23H46	052078-56-5	72



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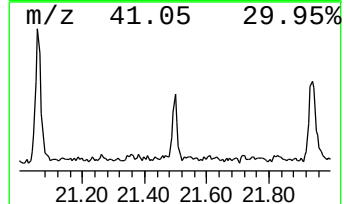
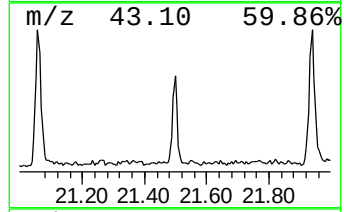
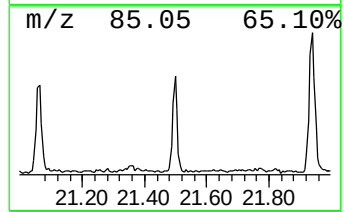
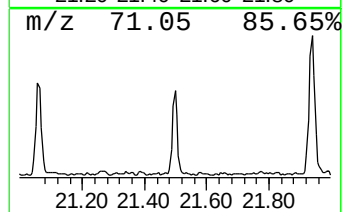
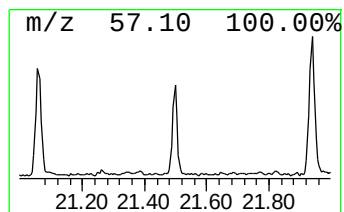
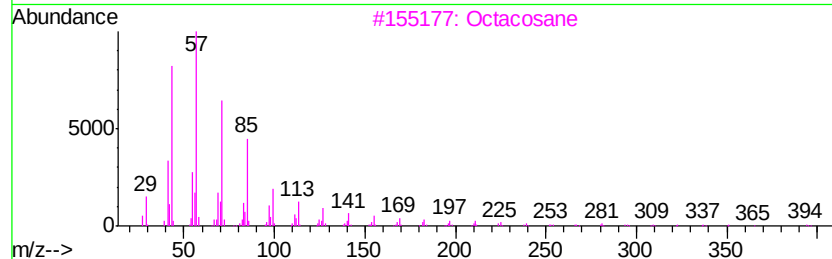
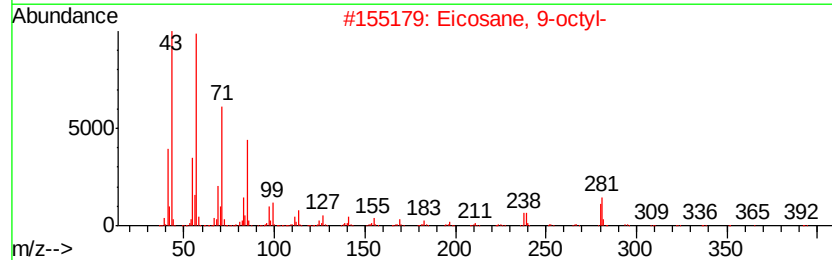
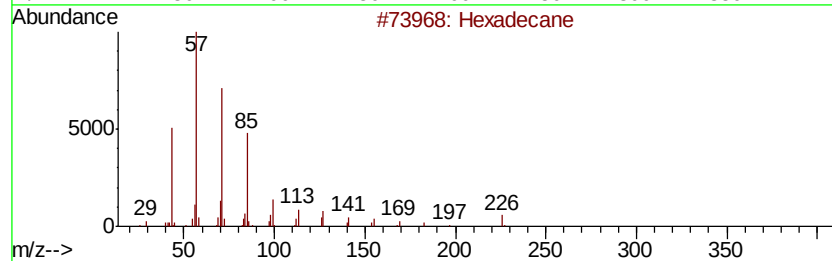
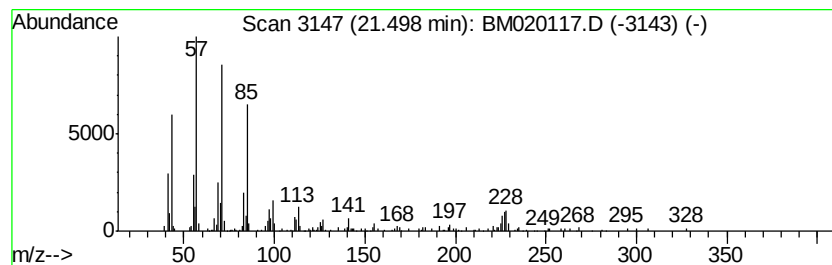
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Hexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.50	2.53 ng	239524	Chrysene-d12	21.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	93
2		Eicosane, 9-octyl-	394	C28H58	013475-77-9	86
3		Octacosane	394	C28H58	000630-02-4	74
4		Heneicosane	296	C21H44	000629-94-7	72
5		Nonacosane	408	C29H60	000630-03-5	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050319\
Data File : BM020117.D
Acq On : 03 May 2019 18:16
Operator : JU/SJ
Sample : K2584-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

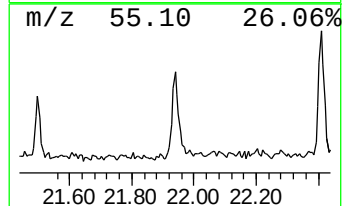
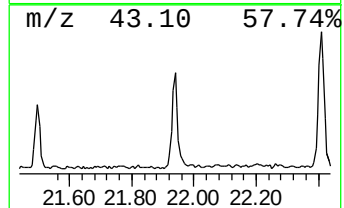
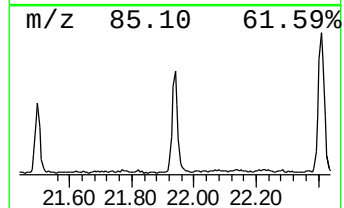
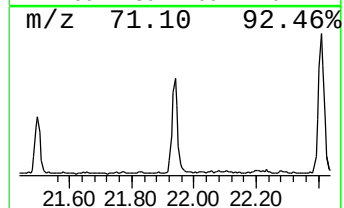
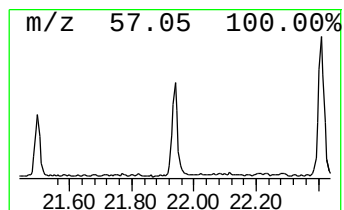
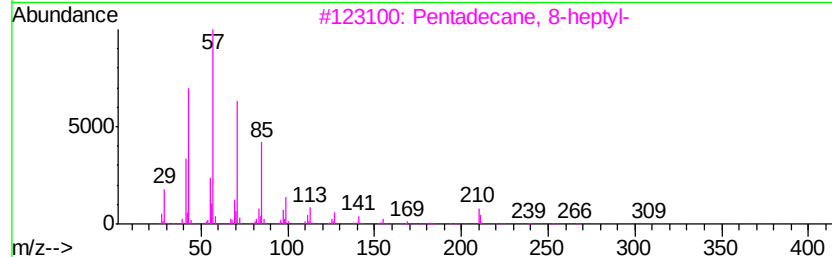
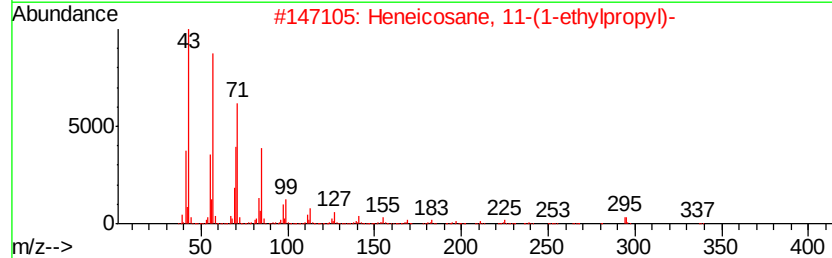
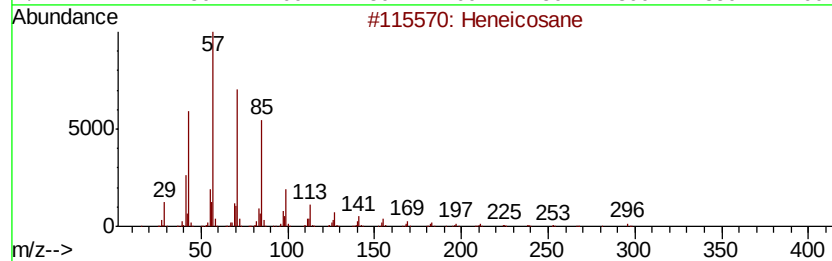
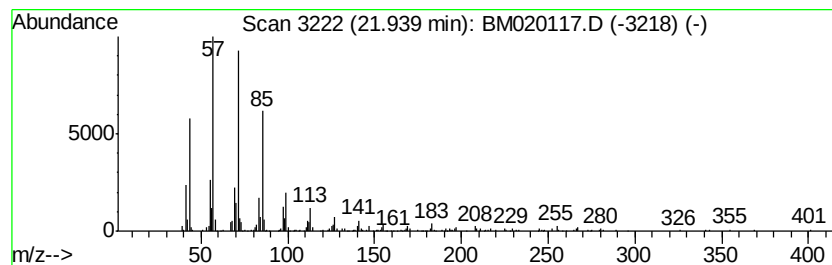
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.94	3.89 ng	367386	Chrysene-d12	21.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	90
3	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90
4	Heptacosane	380	C27H56	000593-49-7	90
5	Octadecane	254	C18H38	000593-45-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050319\
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Sample : K2584-01
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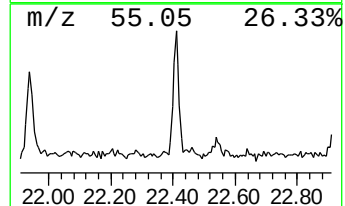
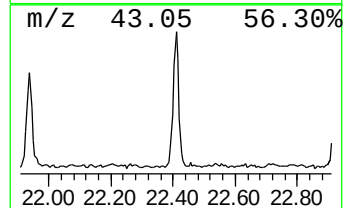
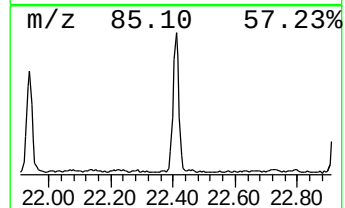
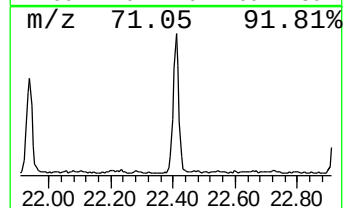
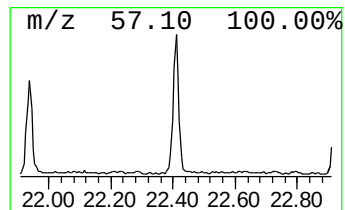
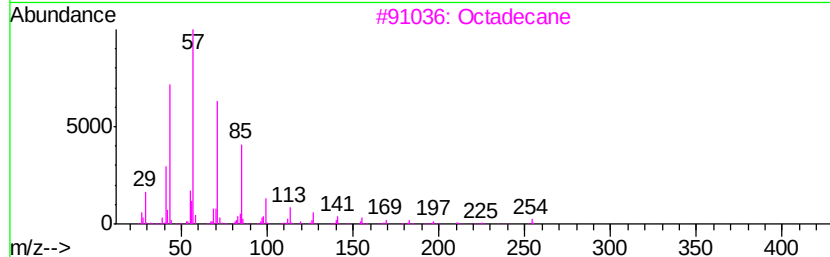
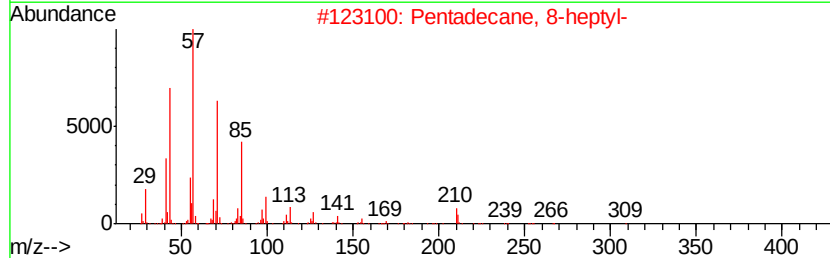
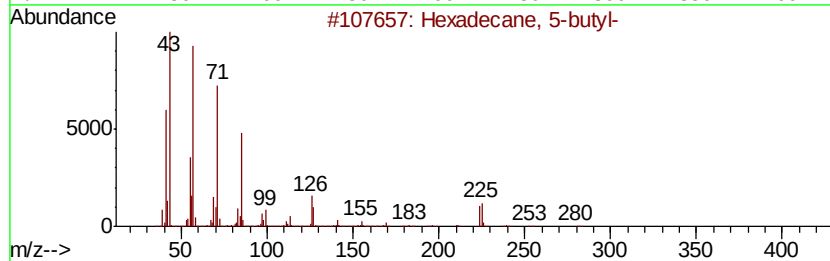
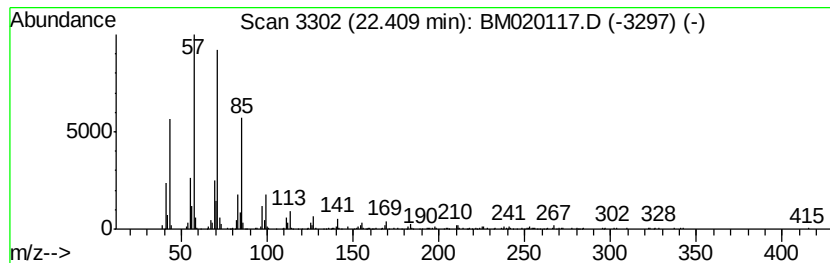
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Hexadecane, 5-butyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.41	5.96 ng	562996	Chrysene-d12	21.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 5-butyl-	282	C20H42	006912-07-8	90
2	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90
3	Octadecane	254	C18H38	000593-45-3	87
4	Heptacosane	380	C27H56	000593-49-7	87
5	Octacosane	394	C28H58	000630-02-4	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050319\
Data File : BM020117.D
Acq On : 03 May 2019 18:16
Operator : JU/SJ
Sample : K2584-01
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Instrument :
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ClientSampleId :
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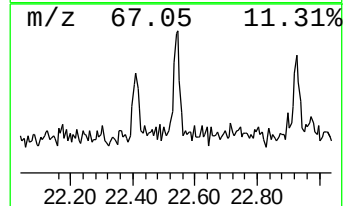
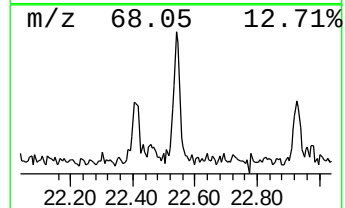
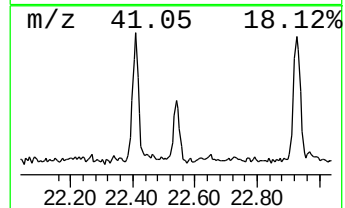
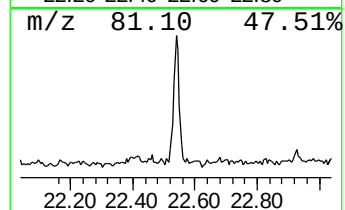
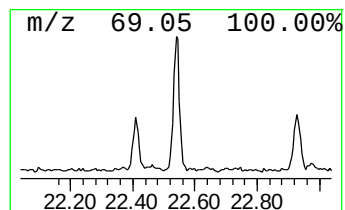
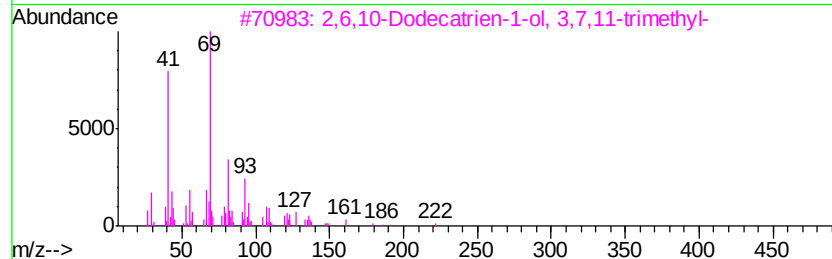
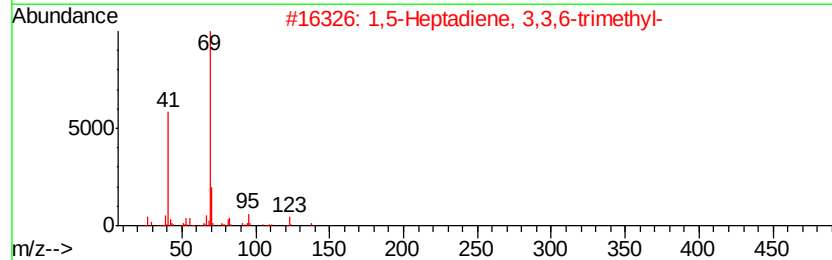
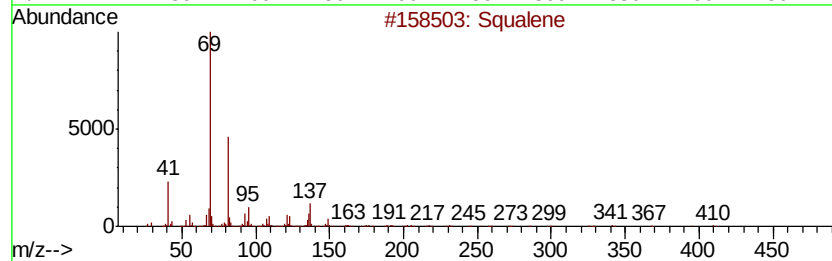
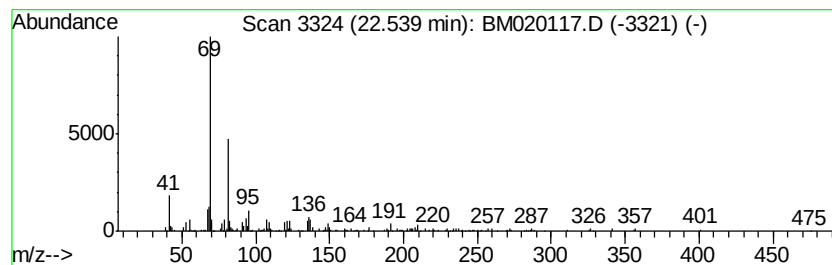
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Squalene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.54	2.14 ng	198063	Perylene-d12	23.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	007683-64-9	74
2		1,5-Heptadiene, 3,3,6-trimethyl-	138	C10H18	035387-63-4	72
3		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	64
4		4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	64
5		2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050319\
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 Acq On : 03 May 2019 18:16
 Operator : JU/SJ
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 ALS Vial : 11 Sample Multiplier: 1

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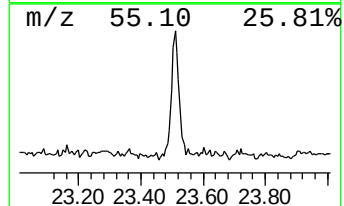
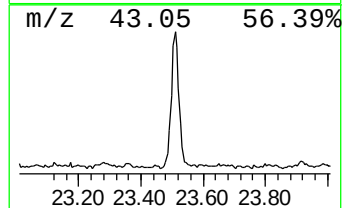
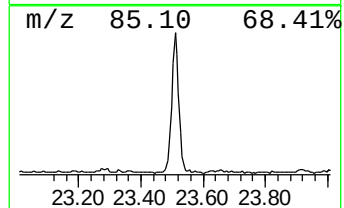
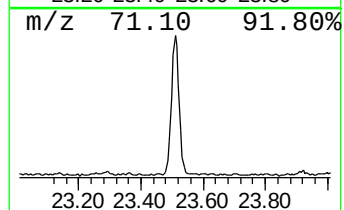
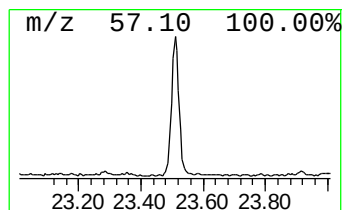
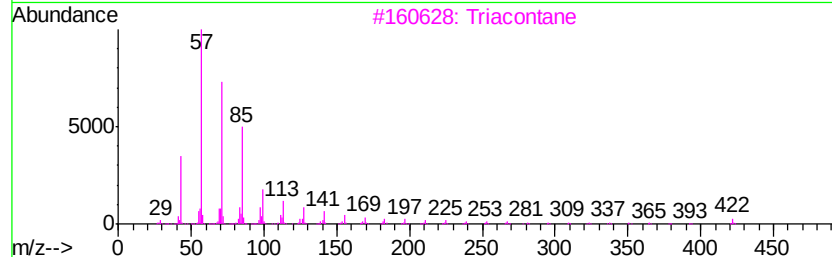
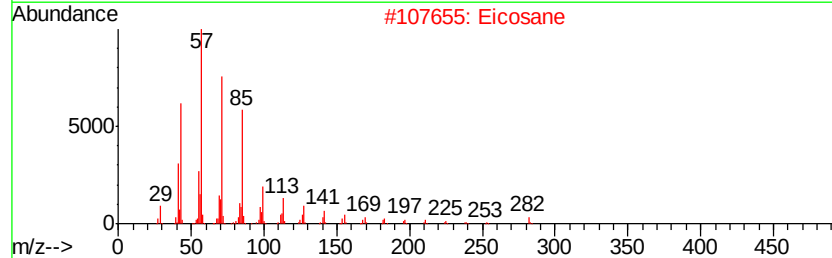
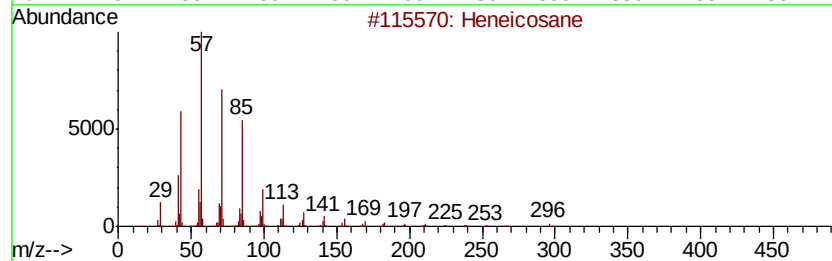
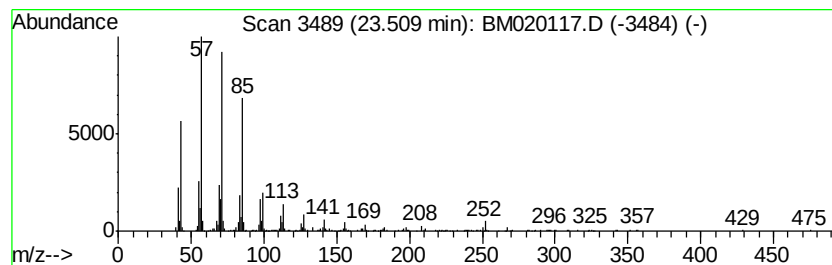
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 11 Triacontane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.51	9.46 ng	874951	Perylene-d12	23.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	90
2		Eicosane	282	C20H42	000112-95-8	87
3		Triacontane	422	C30H62	000638-68-6	86
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	86
5		Hexadecane	226	C16H34	000544-76-3	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050319\
Data File : BM020117.D
Acq On : 03 May 2019 18:16
Operator : JU/SJ
Sample : K2584-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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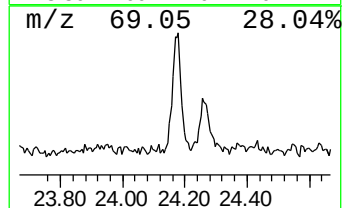
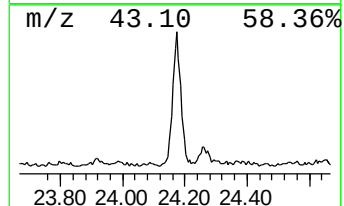
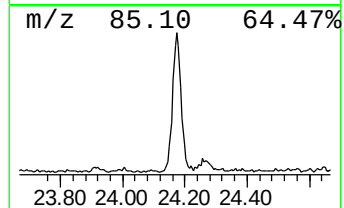
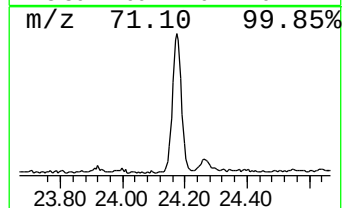
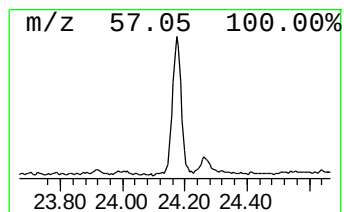
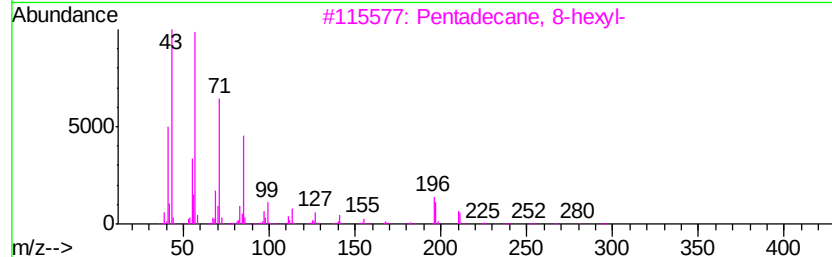
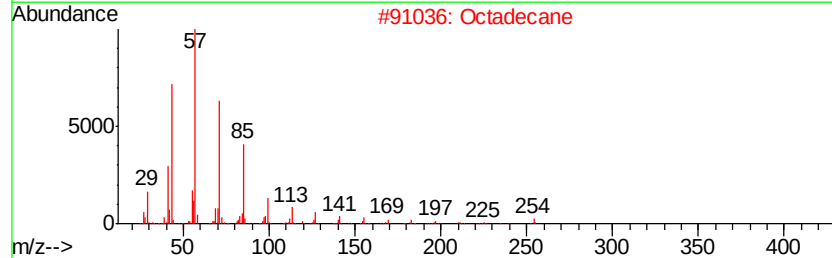
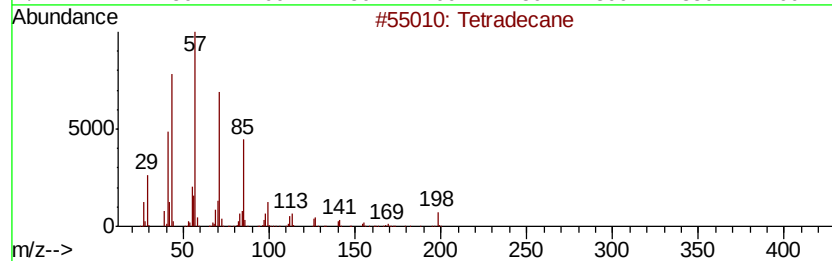
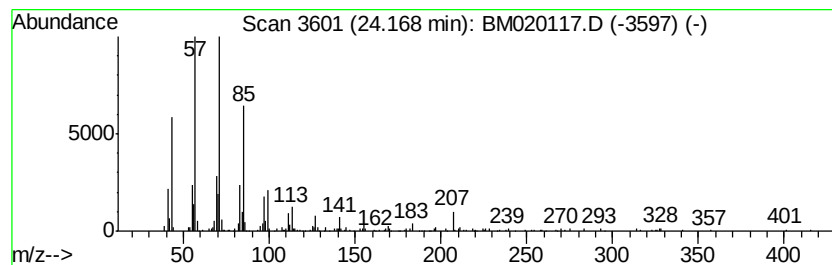
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Tetradecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.17	6.68 ng	618022	Perylene-d12	23.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	87
2			Octadecane	254	C18H38	000593-45-3	87
3			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	87
4			Heptadecane	240	C17H36	000629-78-7	86
5			Octacosane	394	C28H58	000630-02-4	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050319\
Data File : BM020117.D
Acq On : 03 May 2019 18:16
Operator : JU/SJ
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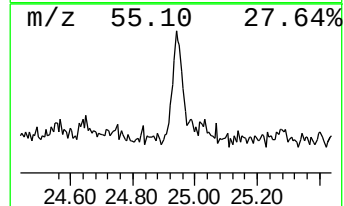
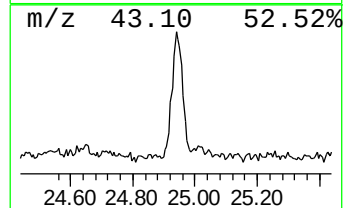
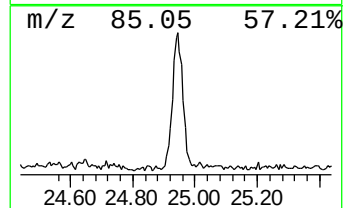
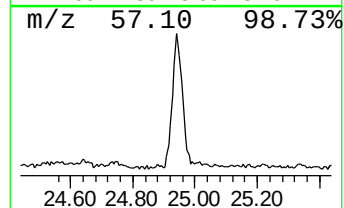
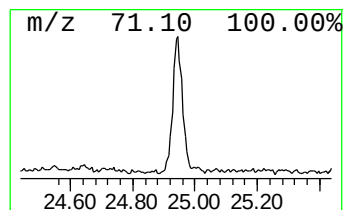
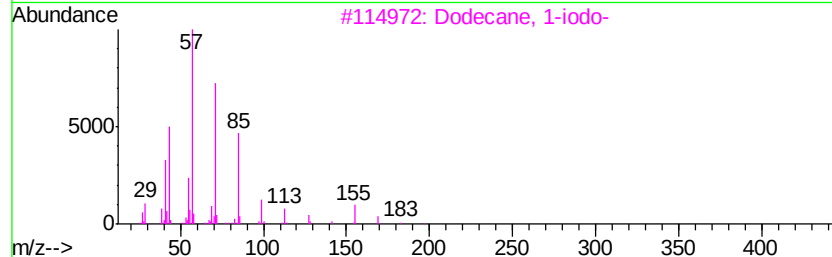
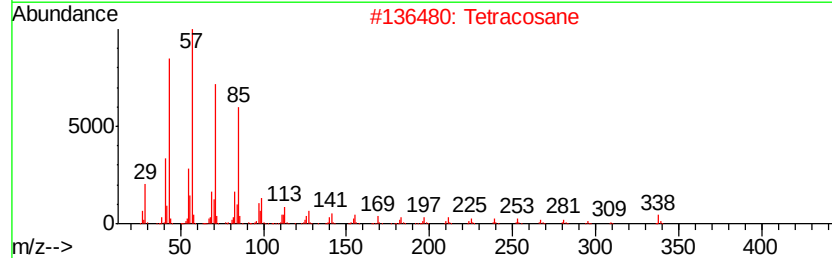
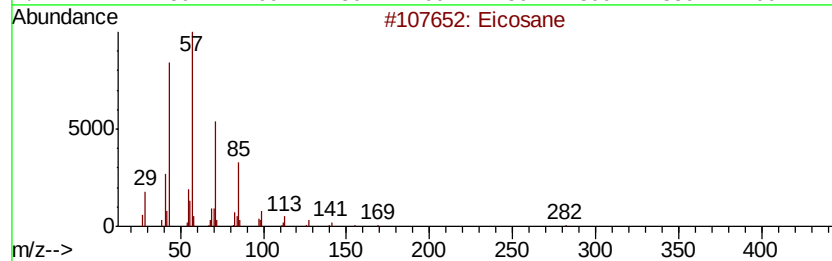
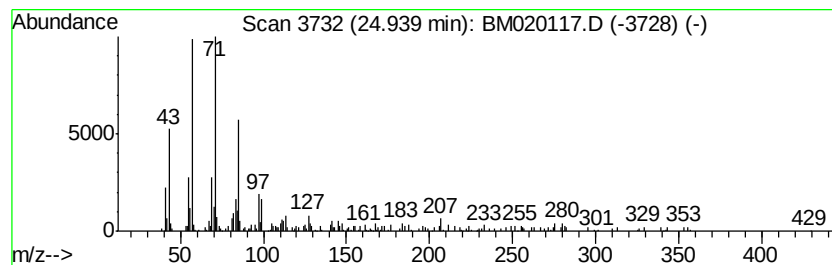
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.94	3.66 ng	338426	Perylene-d12	23.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	93
2		Tetracosane	338	C24H50	000646-31-1	87
3		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	52
5		Octadecane, 3-ethyl-5-(2-ethylbu...	366	C26H54	055282-12-7	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM050319\
 Data File : BM020117.D
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 Sample : K2584-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown3.01	3.01	2.8	ng	104903	1	7.79	743954	20.0
unknown7.32	7.32	140.1	ng	5211400	1	7.79	743954	20.0
n-Hexadecanoic acid	18.06	11.9	ng	823885	4	17.17	1380690	20.0
Octadecanoic acid	19.34	7.0	ng	661020	5	21.36	1890640	20.0
1-Nonadecene	21.06	5.4	ng	509508	5	21.36	1890640	20.0
Hexadecane	21.50	2.5	ng	239524	5	21.36	1890640	20.0
Heneicosane	21.94	3.9	ng	367386	5	21.36	1890640	20.0
Hexadecane, 5-butyl-	22.41	6.0	ng	562996	5	21.36	1890640	20.0
Squalene	22.54	2.1	ng	198063	6	23.64	1850530	20.0
Triacontane	23.51	9.5	ng	874951	6	23.64	1850530	20.0
Tetradecane	24.17	6.7	ng	618022	6	23.64	1850530	20.0
Eicosane	24.94	3.7	ng	338426	6	23.64	1850530	20.0